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AMENDMENT TO THE CLAIMS

The Listing of Claims presented below will replace all prior versions, and listings, of claims in the application:

1. (Currently Amended) A compound having the structure:

$$R_1O$$
 R_2
 R_3
 R_4
 R_5

or pharmaceutically acceptable salt, ester, or salt of such ester; *

wherein R_1 - R_4 R_1 , R_2 and R_4 are each independently hydrogen or an aliphatic, heteroaliphatic, aryl, heteroaryl, alkylaryl or alkylheteroaryl moiety;

R₃ is an aryl moiety having one of the following structures:

 R_5 and R_6 are each independently hydrogen or an aliphatic, heteroaliphatic, aryl, heteroaryl, alkylaryl, or alkylheteroaryl moiety, and wherein R_6 and R_7 , R_5 and R_6 , taken together, may form a cyclic aliphatic, heteroaliphatic, aliphatic(aryl), heteroaliphatic(aryl), aliphatic(heteroaryl) or heteroaliphatic(heteroaryl) moiety, or an aryl or heteroaryl moiety;

wherein each of the foregoing aliphatic and heteroaliphatic moieties may be substituted or unsubstituted, cyclic or acyclic, saturated or unsaturated or linear or branched; and each of the foregoing aryl, heteroaryl, alkylaryl or alkylheteroaryl moieties may be substituted or unsubstituted; and

pharmaceutically acceptable derivatives thereof.

2. (Currently Amended) The compound of claim 1, wherein the compound has the structure (II):

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$$\begin{array}{c|c}
R_1O_{M_{0,1}} & O & R_7 \\
\hline
R_2 & R_4 & R_6 \\
\hline
R_3 & O & R_6 \\
\hline
R_2 & R_6 & R_6
\end{array}$$

$$\begin{array}{c|c}
R_1O_{M_{0,1}} & O & R_6 \\
\hline
R_2 & R_6 & R_6
\end{array}$$

$$\begin{array}{c|c}
R_1O_{M_{0,1}} & O & R_6 \\
\hline
R_3 & R_6 & R_6
\end{array}$$

or pharmaceutically acceptable salt, ester, or salt of such ester;

wherein R_1 , R_4 , R_1 , R_2 and R_4 are each independently hydrogen or an aliphatic, heteroaliphatic, aryl, heteroaryl, alkylaryl or alkylheteroaryl moiety;

R₃ is an aryl moiety having one of the following structures:

 R_5 and R_6 are each independently hydrogen or an aliphatic, heteroaliphatic, aryl, heteroaryl, alkylaryl, or alkylheteroaryl moiety, and wherein R_6 and R_7 , R_5 and R_6 , taken together, may form a cyclic aliphatic, heteroaliphatic, aliphatic(aryl), heteroaliphatic(aryl), aliphatic(heteroaryl) or heteroaliphatic(heteroaryl) moiety, or an aryl or heteroaryl moiety;

wherein each of the foregoing aliphatic and heteroaliphatic moieties may be substituted or unsubstituted, cyclic or acyclic, saturated or unsaturated or linear or branched; and each of the foregoing aryl, heteroaryl, alkylaryl or alkylheteroaryl moieties may be substituted or unsubstituted; and

pharmaceutically acceptable derivatives thereof.

3. (Original) The compound of claim 1, wherein R^1 is hydrogen or an alkyl, heteroalkyl, aryl or heteroaryl moiety substituted with Z, wherein Z is hydrogen, $-(CH_2)_q OR^Z$, $-(CH_2)_q SR^Z$, $-(CH_2)_q N(R^Z)_2$, $-(C=O)R^Z$, $-(C=O)N(R^Z)_2$, or an aliphatic, heteroaliphatic, aryl, heteroaryl, $-(C=O)R^Z$, $-(C=O)R^Z$, -

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USSN 10/649,532 4116711v1 Attorney Docket No.: 2001180-0077 (HU 2060-02 US NATL) (aliphatic)aryl, -(aliphatic)heteroaryl, -(heteroaliphatic)aryl, or -(heteroaliphatic)heteroaryl moiety, wherein q is 0-4, and wherein each occurrence of R^Z is independently hydrogen, a protecting group, a solid support unit, or an aliphatic, heteroaliphatic, aryl, heteroaryl, -(aliphatic)aryl, -(aliphatic)heteroaryl, -(heteroaliphatic)aryl, or -(heteroaliphatic)heteroaryl moiety; wherein each of the foregoing alkyl or heteroalkyl moieties may be substituted or unsubstituted, linear or branched, cyclic or acyclic, saturated or unsaturated; and wherein each of the foregoing aryl, heteroaryl, -(alkyl)aryl, -(alkyl)heteroaryl, -(heteroalkyl)aryl, or -(heteroalkyl)heteroaryl moieties may be substituted or unsubstituted.

- 4. (Original) The compound of claim 3, wherein R¹ is hydrogen, lower alkyl, a substituted or unsubstituted phenyl or -(lower alkyl)phenyl moiety, -(CH₂)_nOR^z, -[(CH₂)_nO]_mR^z, -(CH₂)_n-Ar-(CH₂)_mOR^z; wherein n and m are each independently integers from 1-6, Ar represents a substituted or unsubstituted aryl or heteroaryl moiety, and R^z is independently hydrogen, a protecting group, a solid support unit, or an aliphatic, heteroaliphatic, aryl, heteroaryl, -(aliphatic)aryl, -(aliphatic)heteroaryl, -(heteroaliphatic)aryl, or -(heteroaliphatic)heteroaryl moiety; wherein each of the foregoing alkyl or heteroalkyl moieties may be substituted or unsubstituted, linear or branched, cyclic or acyclic, saturated or unsaturated; and wherein each of the foregoing aryl, heteroaryl, -(alkyl)aryl, -(alkyl)heteroaryl, -(heteroalkyl)aryl, or (heteroalkyl)heteroaryl moieties may be substituted or unsubstituted.
- 5. (Original) The compound of claim 4, wherein R¹ is hydrogen, ethyl, or has one of the structures:

$$R^{2}O$$
 $R^{2}O$
 $R^{2}O$

wherein Rz is as defined in claim 4.

- 6. (Original) The compound of claim 1, wherein R² is hydrogen or an alkyl, heteroalkyl, aryl or heteroaryl moiety substituted with Z, wherein Z is hydrogen, -(CH₂)_qOR^Z, -(CH₂)_qSR^Z, -(CH₂)_qN(R^Z)₂, -(C=O)R^Z, -(C=O)N(R^Z)₂, or an aliphatic, heteroaliphatic, aryl, heteroaryl, -(aliphatic)aryl, -(aliphatic)heteroaryl, -(heteroaliphatic)aryl, or -(heteroaliphatic)heteroaryl moiety, wherein q is 0-4, and wherein each occurrence of R^Z is independently hydrogen, a protecting group, a solid support unit, or an aliphatic, heteroaliphatic, aryl, heteroaryl, -(aliphatic)aryl, -(aliphatic)heteroaryl, -(heteroaliphatic)aryl, or -(heteroaliphatic)heteroaryl moiety; wherein each of the foregoing alkyl or heteroalkyl moieties may be substituted or unsubstituted, linear or branched, cyclic or acyclic, saturated or unsaturated; and wherein each of the foregoing aryl, heteroaryl, -(alkyl)aryl, -(alkyl)heteroaryl, -(heteroalkyl)aryl, or (heteroalkyl)heteroaryl moieties may be substituted or unsubstituted.
- 7. (Original) The compound of claim 6, wherein R² is hydrogen, lower alkyl, a substituted or unsubstituted phenyl or —(lower alkyl)phenyl moiety, —(CH₂)_nOR², —[(CH₂)_nO]_mR², —(CH₂)_n-Ar-(CH₂)_mOR²; wherein n and m are each independently integers from 1-6, Ar represents a substituted or unsubstituted aryl or heteroaryl moiety, and R² is independently hydrogen, a protecting group, a solid support unit, or an aliphatic, heteroaliphatic, aryl, heteroaryl, —(aliphatic)aryl, —(aliphatic)heteroaryl, —(heteroaliphatic)aryl, or —(heteroaliphatic)heteroaryl moiety; wherein each of the foregoing alkyl or heteroalkyl moieties may be substituted or unsubstituted, linear or branched, cyclic or acyclic, saturated or unsaturated; and wherein each of the foregoing aryl, heteroaryl, —(alkyl)aryl, —(alkyl)heteroaryl, —(heteroalkyl)aryl, or (heteroalkyl)heteroaryl moieties may be substituted or unsubstituted.
- 8. (Original) The compound of claim 6, wherein R² is hydrogen or has one of the structures:

wherein R² is as defined in claim 6.

9. (Cancelled)

10. (Cancelled)

- 11. (Original) The compound of claim 1, wherein R⁴ is hydrogen or an alkyl, heteroalkyl, aryl, heteroaryl, -(alkyl)aryl, -(alkyl)heteroaryl, -(heteroalkyl)aryl, or -(heteroalkyl)heteroaryl moiety; wherein each of the foregoing alkyl or heteroalkyl moieties may be substituted or unsubstituted, linear or branched, cyclic or acyclic, saturated or unsaturated; and wherein each of the foregoing aryl, heteroaryl, -(alkyl)aryl, -(alkyl)heteroaryl, -(heteroalkyl)aryl, or (heteroalkyl)heteroaryl moieties may be substituted or unsubstituted.
- 12. (Original) The compound of claim 11, wherein R⁴ is hydrogen alkyl or heteroalkyl.
- 13. (Original) The compound of claim 1, wherein R⁵ and R⁶ are each independently hydrogen or an alkyl, heteroalkyl, aryl, heteroaryl, -(alkyl)aryl, -(alkyl)heteroaryl, (heteroalkyl)aryl, or -(heteroalkyl)heteroaryl moiety; or wherein R⁵ and R⁶, taken together, form a substituted or unsubstituted, saturated or unsaturated cyclic moiety comprising 5-12 carbon atoms, 0-5 oxygen atoms, 0-5 sulfur atoms and 1-5 nitrogen atoms; and wherein each of the foregoing alkyl or heteroalkyl moieties may be substituted or unsubstituted, linear or branched, cyclic or acyclic, saturated or unsaturated; and wherein each of the foregoing aryl, heteroaryl, (alkyl)aryl, -(alkyl)heteroaryl, -(heteroalkyl)aryl, or -(heteroalkyl)heteroaryl moieties may be substituted or unsubstituted.
- 14. (Original) The compound of claim I, wherein -NR⁵R⁶ is one of the following the structures:

15. (Original) The compound of claim 1 having the structure:

16. (Original) The compound of claim 1 having the structure:

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17. (Previously Amended) A compound having the structure:

Claims 18-21: Cancelled

22. (Previously Amended) A compound having the structure:

23. (Currently Amended) A collection of compounds comprising two or more of the compounds of claim 1 or 2 A library of compounds comprising a plurality of library members, wherein at least two library members are a compound of claim 1 or 2.

Claims 24-25: Cancelled

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- 26. (Currently Amended) The collection <u>library</u> of claim 23, wherein the collection <u>library</u> comprises at least 100 compounds.
- 27. (Currently Amended) The collection <u>library</u> of claim 23, wherein the collection <u>library</u> comprises at least 1,000 compounds.
- 28. (Currently Amended) The collection <u>library</u> of claim 23, wherein the collection <u>library</u> comprises at least 2,000 compounds.
- 29. (Currently Amended) The collection <u>library</u> of claim 23, wherein the collection <u>library</u> comprises at least 10, 000 compounds.
- 30. (Currently Amended) A pharmaceutical composition comprising: a compound of any one of claims 1, 2, 5, 8, 10, 14, and 15 22 14, 15-17 and 22; and a pharmaceutically acceptable carrier.

Claims 31-39. Cancelled